In the Claims

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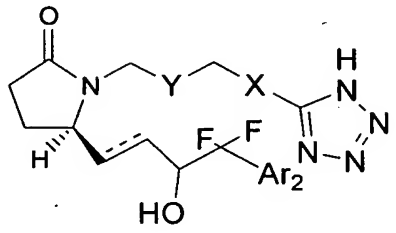
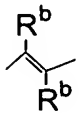
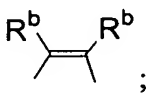
16. Cancel.
17. A compound of structural formula I:



I

or a pharmaceutically acceptable salt, enantiomer, diastereomer, pro drug or mixture thereof, wherein
X is $(CH_2)_n$, O or S;

Y represents $(C(R^b)_2)_n$, triple bond,  or  ;

Ar_2 independently represent $(CH_2)_mC_{6-10}$aryl, $(CH_2)_mC_{5-10}$heteroaryl, $(CH_2)_mC_{3-10}$ heterocycloalkyl, $(CH_2)_mC_{3-8}$ cycloalkyl said cycloalkyl, heterocycloalkyl, aryl or heteroaryl unsubstituted or substituted with 1-3 groups of R_a;

R_a represents C_{1-6} alkoxy, C_{1-6} alkyl, CF_3, nitro, amino, cyano, C_{1-6} alkylamino, or halogen;

R^b independently represents H, halogen, C_{1-6} alkyl, C_{3-6} cylcoalkyl or

--- represents a double or single bond;

n represents 0-4; and

m represents 0-8.

18. The compound according to claim 17 wherein X and Y are $(CH_2)_n$, --- represents a double bond; and Ar_2 is phenyl.

19. The compound according to claim 18 wherein X is $(CH_2)_n$ and n is 1 and Y is $(CH_2)_n$ and n is 3.